Data Limitations and Validation Report
for Lockheed Idaho Technologies
Case No. 93052416, SDG 93052416
Argonne National Laboratory - West
TCL Appendix IX Volatile Organics
Four Aqueous Samples

Validated by:

Michelle L. Allen

Data Validator

Approved by:

Joseph A. Samchuck Data Validation Quality

Assurance Officer

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1.0 INTRODUCTION

The Argonne National Laboratory - West sample set for Case No. 93052416, SDG 93052416 consists of four aqueous samples analyzed for Target Compound List (TCL) Appendix IX volatile organic compounds. The analyses were performed by Biospherics, Inc. using the protocols outlined in the "Analytical Laboratory Services for Environmental Groundwater Samples for the Argonne National Laboratory - West" Statement of Work (SOW). A total of 184 TCL sample data points were reported in this analytical data set.

The analytical data from these analyses were reviewed by HALLIBURTON NUS Corporation personnel as Level A in accordance with ERP Standard Operating Procedure SMO-SOP-12.1.3.

2.0 QUALITY CONTROL SUMMARY

The data were evaluated based on the following parameters:

Data Completeness

- Holding Times
 GC/MS Tuning and Mass Calibration
 Initial and Continuing Calibrations
 Blank Analyses
- Surrogate Spike Recoveries
- Matrix Spike/Matrix Spike Duplicate Results Blank Spike Results
- * Internal Standards Performance
- System Performance and Detection Limits Laboratory Performance
- * Compound Quantitation

The asterisk indicates that all quality control criteria were met for this parameter. Problem areas affecting data usability are discussed in Section 4.0 of this report. A Glossary of Data Validation Flags which defines the validation qualifiers applied on a sample-specific basis is presented in Section 6.0.

3.0 DATA COMPLETENESS

The data presented in Case No. 93052416, SDG 93052416 consists of TCL Appendix IX volatile organic results for four (4) aqueous sample as follows:

93052416-1 (EBR II NO1) 93052416-2 (EBR II NO2) 93052416-3 (MW-11) 93052416-4 (TRIP BLANK)

The data package was missing the bromofluorobenzene (BFB) tune (Form V) corresponding to the initial calibration, the initial calibration Form VI containing target compound Percent Relative Standard Deviations (%RSDs) and Relative Response Factors (RRFs), and the associated initial calibration raw data. In addition, the Blank Spike (BS) analysis Form III, and the laboratory method blank Form I and associated raw data were not included in this data package. Hence, the data in this SDG could not be evaluated for these parameters. It should be noted that the presentation and documentation of the data package deliverables were extremely poor. The data package does not conform to a Level A deliverable.

4.0 SUMMARY OF DATA USABILITY

The Matrix Spike/Matrix Spike Duplicate (MS/MSD) samples were analyzed outside the 12 hour BFB tune. No action was taken since these are quality control samples.

Continuing calibration Percent Differences (%Ds) for chloromethane, bromomethane, chloroethane, methylene chloride, acetone, carbon disulfide, and xylenes (total) exceeded 75%. The nondetected results reported for these compounds in the associated environmental samples were qualified as rejected, "R".

Some continuing calibration %Ds for 1,1-dichloroethene, 2-butanone, cis-1,3-dichloropropene, trans-1,3-dichloropropene, and 1,1,2,2-tetrachloroethane were greater than the 25% quality control limit. No actions were necessary since no positive results were reported for these compounds in the affected samples and the nondetects were not compromised.

The MS/MSD analyses yielded Percent Recoveries (%Rs) for 1,1-dichloroethene and benzene were above the upper quality control limits. No qualifications were necessary since no positive results were reported for these compounds in the unspiked sample.

The Appendix IX volatile compounds allyl chloride, methacrylonitrile, and propanenitrile were included in the continuing calibration 50 ppb standard. Hence, these compounds were qualified as rejected, "R",on the environmental sample Form Is.

Annotated laboratory Form I data summary reports showing the data and relevant qualifier flags applied are presented in Appendix A of this report. Copies of the unqualified data summary reports as reported by the laboratory are provided in the attached Appendix B. The attached Appendix C includes documentation to support the findings discussed in this report.

A sample-specific summary of the data validation flags applied is depicted in Table 1, appearing on the following page. The qualifier flags used as a result of the validation process are defined in Section 6.0 (Glossary of Data Validation Flags) of this report. Details regarding the application of the validation qualifiers are discussed in the remainder of this section.

4.1 GC/MS Tuning

The MS/MSD samples were analyzed outside the 12 hour BFB tune criterion. No actions were necessary since these are quality control samples.

4.2 Calibrations

The continuing calibration performed on instrument GC/MS#1 (05/28/93) contained the following continuing calibration Percent Differences (%Ds) which exceeded 75%.

| Compound | <u>%D</u> |
|--------------------|-----------|
| chloromethane | 70.6 |
| | 79.6 |
| bromomethane | 93.0 |
| chloroethane | 110.7 |
| methylene chloride | 126.8 |
| acetone | 117.0 |
| carbon disulfide | 76.8 |
| xylenes (total) | 132.5 |

TABLE 1 Lockheed Idaho Technologies Argonne National Laboratory - West Case No. 93052416, SDG 93052416 TCL Appendix IX Volatile Organics

| 93052416-1 (EBR II NO1) R ^{1,2} 93052416-2 (EBR II NO2) R ^{1,2} 93052416-3 (MW-11) R ^{1,2} 93052416-4 (TRIP BLANK) R ^{1,2} | Sample No. | Qualifier Flags |
|---|---|---------------------------------------|
| | 93052416-2 (EBR II NO 93052416-3 (MW-11) | D2) R ^{1,2} R ^{1,2} |

Affected Samples:

ΑII

The nondetected results reported for these compounds in the associated environmental samples were qualified as rejected, "R".

The continuing calibration performed on instrument GC/MS#1 (05/28/93) contained the following continuing calibration Percent Differences (%Ds) which failed to meet the 25% quality control criteria.

| Compound | <u>%D</u> |
|---|--------------------------------------|
| 1,1-dichloroethene 2-butanone cis-1,3-dichloropropene trans-1,3-dichloropropene 1,1,2,2-tetrachloroethane | 28.6 34.1 45.1 41.1 43.3 |
| Affected Sample: All | 40.0 |

Only nondetected results were reported for these compounds in the affected samples and these nondetects were not compromised.

4.3 Matrix Spike/Matrix Spike Duplicate

The %Rs for 1,1-dichloroethene and benzene in the MS/MSD analyses performed on sample 93052416-4 (TRIP BLANK) were high. No qualifications were necessary since no positive results were reported for these compounds in the unspiked sample.

5.0 SUMMARY OF LABORATORY PERFORMANCE

The data associated with the initial calibration was missing from this SDG. The BS analysis Form III was not included. The field identification was not used to name the environmental samples, the sample location ("EBR II NO1, EBR II NO2, MW-11, and TRIP BLANK") were used. Hence, the data validator used both the Biospherics laboratory IDs, 93052416-1, -2, -3, and -4, and the field IDs. The data validator was not positive that the correct Chain-of-Custody (C-O-C) was referenced for the data validation as a result of the incorrect field IDs used. Several compounds were rejected as a result of %Ds exceedances >75%. It should be noted that several forms had incomplete and/or incorrect times and dates, the internal standard Form VIII had the wrong continuing calibration areas/retention times, and the sample Form Is did not contain laboratory qualifiers. The data validator manually corrected these errors. A laboratory method blank was not included in the data package. Xylenes (total) and the individual isomers (m,p-xylene and o-xylene) were both reported on the sample Form Is. Validation action was taken only on xylenes (total). The laboratory failed to standardize for three target compounds, hence, these were rejected on the sample Form Is. Several continuing calibration %Ds were noncompliant.

6.0 GLOSSARY OF DATA VALIDATION FLAGS

The following data validation flags were applied to the sample data for reasons detailed previously in this report:

R¹ - Reject, "R", nondetected results for chloromethane, bromomethane, chloroethane, methylene chloride, acetone, carbon disulfide, and xylenes (total) as a result of continuing calibration %Ds > 75%.

 Reject, "R", nondetected results for allyl chloride, methacrylonitrile, and propanenitrile as a result of the laboratory's failure to standardize for these compounds.

7.0 REFERENCES

The data referenced in this report were validated in accordance with the protocols outlined in ERP Standard Operating Procedure SMO-SOP-12.1.3 as presented in ERP-SOW-37. In addition, details stipulating laboratory procedures as outlined in "Analytical Laboratory Services for Environmental Groundwater Samples for the Argonne National Laboratory - West" SOW were referenced.

APPENDIX A QUALIFIEDLABORATORYRESULTS

| Lab Name: BI | ospheri | CS INCORP | ORATED | Contract: | ARGONNE | | BR II N | 01 | 1 1 (| |
|--|--|---------------------------------------|-------------|--|---|-------|---------|-------------|-------------|----------------|
| Lab Code: EB | RIINO1 | Case No. | 93052416 | SAS No.: | SDG | No. : | | — <u></u> - | | |
| Matrix: (soil, | /water) | WATER | | | Lab Sample ID: | 930 | 52416-1 | | | |
| Sample wt/vol: | • . | | (g/mL) | | Lab File ID: | 824 | 0 | | | |
| Level: (low/ma | ed) | LOW | | | Date Received: | 05/ | 21/93 | | | |
| Moisture: no | ot dec. | · · · · · · · · · · · · · · · · · · · | | | Date Analyzed: | 05/ | 28/93 | | | |
| GC Column: | · <u> </u> | ID: | | (mm) | Dilution Factor: | 1 | | | | |
| Soil Extract V | /olume: | | | (uL) | Soil Aliquot Volum | ne: | | (| uL | |
| CAS | NO. | | (| COMPOUND | CONCENTRATION (ug/L or ug/ | | | | Q | |
| 74 75 75 75 75 75 75 75 75 76 77 70 79 71 100 79 71 100 79 71 100 79 112 | -83-9- -01-4- -00-3- -09-2- -64-1- -15-0- -35-4- -34-3- -59-0- -66-3- -7-06-2- -93-3- -27-4- -87-5- -01-6- 4-48-1- -01-6- 4-48-1- -00-5- -43-2- 061-02- -25-2- 8-10-1- 1-78-6 7-18-4- -34-5- 8-88-3- 8-90-7- 0-42-5- | | | Bromometha Vinyl Chlo Chloroetha Acthylene Acetone Carbon Dis 1,1-Dichlo 1,2-Dichlo Chloroform 1,2-Dichlo Carbon Tet Bromodichlo 1,2-Dichlo Carbon Tet Car | ne ride ride ne Chloride clfide roethene roethane roethane roethane hloroethane rachloride coromethane ropropane chloropropene thene coromethane nloroethane rothoroethane coromethane | 2 | | 10 | | 2 |
| I . | | | _ | ylene (to | tal) | | / | 51 V | Z · | R |

VOLATILE ORGANIC ANALYSIS DATA SHEET

| Lab Name: BIOSPHERICS INCORPO | ORATED | Contract: | ARGONNE | EBR II NO1 |
|-------------------------------|----------|---|--|--|
| Lab Code: EBRII NO1 Case No. | 93052416 | SAS No.: | SDG No. | · |
| Matrix: (soil/water) WATER | 1 | • | Lab Sample ID: | 93052516-1 |
| Sample wt/vol: | (g/mL) | | Lab File ID: | 8240 |
| Level: (low/med) LOW | | | Date Received: | 05/21/93 |
| % Moisture: not dec | | | Date Analyzed: | 05/28/93 |
| GC Column: ID: | | (mm) | Dilution Factor: | |
| Soil Extract Volume: | | (uL) | Soil Aliquot Volume: | (uL |
| oil Extrac CAS NO. | | COMPOUND | CONCENTRATION UNI (ug/L or ug/Kg) | |
| 97-63-2 | | Acrolein Vinyl Acet Acrylonitr 2-Chloroet M & P Xyle O Xylene Methyl ter Allyl Chlo Mothacrylo Propencit Lodomethan | tate tile thyl Vinyl Ether the the tt-Butyl Ether tride v. | MA 5 4 2 27 16 50 10 10 10 10 50 50 50 50 50 50 50 50 50 50 50 50 50 |

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FORM I VOA

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| Lab Name: | BIOSPHERI | CS INCORP | ORATED | Contract: | ARGONNE | EBR II NO2 | |
|-------------|-----------------------|------------|----------------|-----------------------|--------------------|------------|--------------|
| Lab Code: | EBRIINO2 | Case No. | 93052416 | SAS No.: | SDG | Жо. : | —— |
| Matrix: (s | oil/water) | WATER | _ | | Lab Sample ID: | 93052416-2 | |
| Sample wt/ | vol: | <u> </u> | (g/mL) | | Lab File ID: | 8240 | |
| Level: (lo | w/med) | LOW | <u>-</u> | | Date Received: | 05/21/93 | |
| % Moisture: | : not dec. | | • | | Date Analyzed: | 05/28/93 | |
| GC Column: | | ID: | | (mm) | Dilution Factor: | 1 | |
| Soil Extra | ct Volume: | | | (uL) | Soil Aliquot Volu | me:(| uL |
| | | | | | CONCENTRATION | INTTQ. | |
| | CAS NO. | | | COMPOUND | (ug/L or ug | | Q |
| | 74-87-3- | | | Chlorometh | hane | MA 10-11 | |
| | 74-83-9- | | | Bromometha | ne | 22196 10 | 耳と |
| | 75-01-4- | | | Vinyl Chlo | oride | 10, 6 | T. |
| i | 75-00-3- | | | Chloroeth | ine | 10 4 | FR |
| 1 | 1 /5-09-2- | | · : | Methylene | Chloride | 5+6 | ER |
| | 75-15-0- | | | Acetone Carbon Dia | | 10 | ER |
| | 75-15-0- | | | 1,1-Dichlo | antide | | FIR |
| | 75-34-3- | | . | 1,1-Dichle | oroethane | 51.4 | ــنــ |
| | 540-59-0- | | | 1.2-Dichle | proethene (total) | 51 | |
| | 67-66-3- | | - - | Chloroform | n | 51 | ب_ |
| ! | 107-06-2- | | | 1,2-Dichlo | roethane | 5: | — |
| | 78-93-3- | | : | 2-Butanone | • | 101 | |
| i | 71-55-6- | | - - | 1,1,1-Tric | chloroethane | 5; | <u> </u> |
| ! | 56-23-5- | | · (| Carbon Tet | rachloride | 51 | |
| į | 75-27-4- | | · ! | Bromodichl | oromethane | 5 | |
| ; | 78-87-5- | | : | 1,2-Dichlo | ropropane | 5! | |
| į | | -5 | | cis-1,3-Di | chloropropene | 5, | \Box |
| į | /9-01-6- 124-48-1- | | | Trichloroe | | 5! | |
| ¦ | 79-00-5- | | · | | oromethane | 51 | |
| | 71-43-2- | | | 1,1,2-Tric Benzene | chloroethane | 5. | |
| 1 | 10061-02- | | • | | . D.(=b.) ======= | 51 | |
| l L | 75-25-2- | | | Bromoform | -Dichloropropene | 5 | |
| į | 108-10-1- | - - | : | 4-Methul-2 | ?-Pentanone | 51 | |
| : | 591-78-6 | | · : | 2-Hexanone | · renrations | 10, | |
| į | 127-18-4- | | | Tetrachlor | coethene | 10: | - ' |
| i | | | | 1,1,2,2-Te | trachloroethane | 5; | <u>—</u> |
| ! | 108-88-3- | | | Toluene | | 51 | — |
| į | 108-90-7- | | | Chlorobenz | zene | 51 | |
| 1 | | | : | Ethylbenze | | 5 | - |
| ! | 100-42-5- | | | Styrene | | 5 | |
| i | 1330-20-7 | 7 | · : | Xylene (to | otal) | L Sik | |
| ļ. | | | | | | | <u>~</u> ;`` |

| Lab Name: 1 | BIOSPHERIC | S INCORP | ORATED | _Contract: | ARGONNE | EBR | R II NO |)2 | |
|------------------------------|-----------------------|-----------|-------------|--------------|----------------------------|----------|----------------|------------|--------|
| Lab Code: 1 | BRII NO2 | Case No. | 93052416 | _ SAS No.: | SDG | No. : | | | |
| Matrix: (soi ab Code: | ll/water)_ | WATER | | • | Lab Sample ID: | 9305 | 2416-2 | | |
| Sample wt/vo atrix: (soi) | | | (g/mL) | | Lab File ID: | 8240 | | | |
| Level: (low/ ample wt/vol | med) | LOW | | | Date Received: | 05/2 | 1/93 | _ | |
| Moisture: | | | | | Date Analyzed: | 05/2 | 8/93 | | |
| GC Column: Moisture: n | | ID: | | (mm) | Dilution Factor: | | | _ | |
| Soil Extract C Column: | Volume: | - | | (uL) | Soil Aliquot Volum | ne: | | _ (ı | υL |
| oil Extrac C | AS NO. | | | COMPOUND | CONCENTRATION (ug/L or ug/ | | | _ ' | Q |
| | 97-63-2 107-02-8- | · | | Ethyl Met | hacrylate | M | A | 5 6 | |
| ! : | 108-05-4- | | | Vinul Acet | tate | <u> </u> | | | \Box |
| (10 | 07-13-1 | · | | Acrul and to | rila | <u> </u> | | 0 | H |
| | 110-75-8- | | | 2-Chloroet | hel Vinul Pehan | | | 0 | |
| | 700-30-3- | | | Μ ፋ Þ Υυία | ine | | 1 | 0 | \neg |
| | 95-47-6 1634-04-4- | · | | O Xylene | | | | 5! | \Box |
| | 17-05-1 | · | | Metnyi te: | t-Butyl Ether | | | 51 | 三 |
| 11 | 26-98-7~ - | · | | Manhaanal | ride. | | | 6 / | \Box |
| 1 1 | 107-12-0- | | | Recommend | materia. | ļ | | OI A | |
| ļ 7 | 74-88-4 | | | Iodomether | | <u> </u> | | O R | |
| 1 8 | 90-62-6 | | 1 | Methyl Met | chacrylate | <u> </u> | √ 2 | 0; N | L |
| <u></u> | | | | | | <u>.</u> | | _!_ | |

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FORM I VOA

| | | | | | į | |
|-------------|------------|-------------|--------------|----------------|--------------------------|---------------|
| Lab Name: | BIOSPHERI | CS INCORP | ORATED | Contract: | ARGONNE | ₩-11 |
| Lab Code: | MW-11 | Case No. | 93052416 | SAS No.: | SDG | Жо.: |
| Matrix: (s | oil/water) | WATER | | | Lab Sample ID: | 93052416-3 |
| Sample wt/ | vol: | | (g/mL) | | Lab File ID: | 8240 |
| Level: {lo | w/med) | LOW | | | Date Received: | 05/21/93 |
| % Moisture | not dec. | | | | Date Analyzed: | 05/28/93 |
| GC Column: | | ID: | | (mm) | Dilution Factor: | 1 |
| Soil Extrac | ct Volume: | | | (uL) | Soil Aliquot Volu | me:(uL |
| | | | | | CONCENTRATION | UNITS: |
| | CAS NO. | | | COMPOUND | (ug/L or ug | /Kg) UG/L Q |
| | 1 74 07 3 | | | Chlorometh | | |
| | 1 74-07-3- | | | Bromometha | nane | MA 10 UT |
| ļ | 75-01-4- | | | Vinyl Chlo | ine setda | 2/27/76 10 UT |
| | 75-00-3- | | | Chloroetha | oride | 10 U |
| ! | 75-09-2- | | | Methylene | Chlorida | 10,45 |
| | 67-64-1- | | | Acetone | CHIOITE | 10117 |
| | 75-15-0- | | | Carbon Dis |) #1.da | 10,14 |
| i | 75-35-4- | | . - | 1,1-Dichlo | artine | 51 117 |
| | 75-34-3- | | | 1,1-Dichic | roethana | 5 K |
| į | 540-59-0- | | | 1, 2-Dichie | proethene (total) | 51 U |
| | 67-66-3- | | . - | Chloroform | - (total) | <u> </u> |
| | 107-06-2- | | | 1,2-Dichle | roethane | 51 U |
| | 78-93-3- | | | 2-Butanone | toechane | 51 |
| | 71-55-6- | | | 1 1 1 mm = 4 a | : :hloroethane | 101 |
| | 56-23-5- | | · - | Carbon Tet | rachloride | 5 |
| ! | 75-27-4- | | - - | Bromodichl | oromethane | 51 |
| | 78-87-5- | | . - | 1.2-Diahla | coromethane copropane | 5 |
| ! | 10061-01- | -5 | | cis-1.3-ni | chloropropene | 5 |
| i | 79-01-6- | | | Trichloroe | thene | 5 |
| 1 | 124-48-1- | _ _ | - - - | Dibromochi | oromethane | 5 |
| į | 79-00-5- | | | | chloroethane | 51 -1 |
| 3 | | | | Benzene | | 5 |
| į | 10061-02- | -6 | | trans-1.3- | Dichloropropene | 5! -! |
| i | 75-25-2- | | | Bromoform | - | 3 |
| ! | 108-10-1- | · | | 4-Methvl-2 | -Pentanone | 10 |
| i | 1 591-78-6 | | | 2-Hexanone | | |
| ! | 127-18-4- | | | Tetrachlor | coethene | 5: |
| i | 79-34-5- | | | 1.1.2.2-Te | trachloroethane | 51 |
| ļ | 108-88-3- | | | Toluene | | 31 - 31 |
| į | 108-90-7- | | | Chlorobenz | ene | 51 |
| | 100-41-4- | | | Ethylbenze | ene | 5 / |
| į | 100-42-5- | | | Styrene | | 31 1 |
| | 1330-20-7 | 1 | | Xylene (to | otal) | -5.4m |
| | | | | - | | |

| Lab Name: 1 | Biospheri | CS INCORP | ORATED | _Contract: | ARGONNE | MW-11 | |
|--------------|--|-----------|----------|---|--|---|-----|
| Lab Code: 1 | MV-11 | Case No. | 93052416 | SAS No.: | SDG N | lo.: | |
| Matrix: (soi | il/water) | WATER | | • | Lab Sample ID: | 93052416-3 | |
| Sample wt/vo | ol: | | (g/mL) | | Lab File ID: | 8240 | |
| Level: (low/ | /med) | LOW | | | Date Received: | 05/21/93 | |
| Moisture: | not dec. | | | | Date Analyzed: | 05/28/93 | |
| GC Column: _ | | ID: | | (mm) | Dilution Factor: | | |
| Soil Extract | Volume: | | | (uL) | Soil Aliquot Volume | B:(| (uL |
| oil Extrac C | AS NO. | | | COMPOUND | CONCENTRATION (ug/L or ug/l | 10 _ L | Q |
| 1 | 107-02-8- 108-05-4- 07-13-1- 110-75-8- 108-38-3- 95-47-6- 1634-04-4 17-05-1- 26-98-7- 107-12-0- 74-88-4- | | | Vinyl Acet Acrylonit; 2-Chloroet M & P Xyle O Xylene Nethyl ter Allyl Ghlo Methacrylo Respensit | tate tile thyl Vinyl Ether the t-Butyl Ether tride | 50 21214, 50 10 50 10 10 5 5 50 50 50 50 50 | 4 |

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FORM I VOA

| Lab Name: | BIOSPHERI | CS INCORE | PORATED | Contract: | ARGONNE | TRIP BLANK | 1 |
|------------|--|-----------|----------|--|--|-----------------|-----|
| Lab Code: | BLANK | Case No. | 93052416 | SAS No.: | SDG | No.: | |
| Matrix: (s | oil/water) | WATER | _ | | Lab Sample ID: | 93052416-4 | |
| Sample wt/ | vol: | | (g/mL) | | Lab File ID: | 8240 | |
| Level: (lo | w/med) | LOW | _ | | Date Received: | 05/21/93 | |
| % Moisture | not dec. | | _ | | Date Analyzed: | 05/28/93 | |
| GC Column: | | ID: | | (mm) | Dilution Factor: | 1 | |
| Soil Extra | ct Volume: | | | (uL) | Soil Aliquot Volu | me: | (uL |
| | a.a. wa | | | | CONCENTRATION | | |
| | CAS NO. | | | COMPOUND | (ug/L or ug | /Kg) UG/L | Q |
| | 74-83-9- 75-01-4- 75-00-3- 75-09-2- 67-64-1- 75-15-0- 75-35-4- 75-34-3- 540-59-0- 67-66-3- 107-06-2- 78-93-3- 71-55-6- 56-23-5- 75-27-4- 78-87-5- 10061-01- 79-01-6- 124-48-1- 79-00-5- 71-43-2- 10061-02- 75-25-2- 108-10-1- 591-78-6 127-18-4- | | | Carbon Dis 1,1-Dichlo 1,1-Dichlo 1,2-Dichlo Chloroform 1,2-Dichlo 2-Butanone 1,1,1-Tric Carbon Tet Bromodichl 1,2-Dichlo cis-1,3-Di Trichloroe Dibromochl 1,1,2-Tric Benzene trans-1,3- Bromoform 4-Methyl-2 2-Hexanone Tetrachlor | ene Oride ane Chloride sulfide proethene proethene proethane crachloride crachloride coromethane crachloropropene chloropropene | 101 51 51 | |
| | 100-41-4- 100-42-5- 1330-20-1 | | | Ethylbenze Styrene Xylene (to | ene | 51 51 | |

| Lab Name: BIOSPHERICS INCORP | ORATED | _Contract: | ARGONNE | TRIP BLANK |
|------------------------------|----------|--|--|--|
| Lab Code: BLANK Case No. | 93052416 | SAS No.: | SDG No | ·: |
| Matrix: (soil/water) WATER | | • | Lab Sample ID: | 93052416-4 |
| Sample wt/vol: | (g/mL) | | Lab File ID: | 8240 |
| Level: (low/med) LOW | , | | Date Received: | 05/21/93 |
| % Moisture: not dec. | | | Date Analyzed: | 05/28/93 |
| GC Column: ID: | | (mm) | Dilution Factor: | |
| Soil Extract Volume: | | (uL) | Soil Aliquot Volume: | |
| oil Extrac CAS NO. | | COMPOUND | CONCENTRATION UN (ug/L or ug/Kg | |
| 97-63-2 | | Acrolein Vinyl Acet Acrylonitr 2-Chloroet 4 & P Xyle 5 Xylene 4ethyl ter 41yl-Chlo 6etharylo | ate ile hyl Vinyl Ether ne t-Butyl Ether side- nitrile | 5 LL 2/27/4/50 10 10 10 10 5 5 5 60 R 60 R 60 R 50 U |

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FORM I VOA